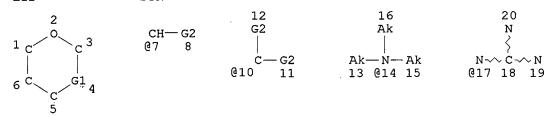
=> fil reg FILE 'REGISTRY' ENTERED AT 09:20:07 ON 01 JUL 1999 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 1999 American Chemical Society (ACS)

STRUCTURE FILE UPDATES: 01 JUL 99 HIGHEST RN 226917-44-8 DICTIONARY FILE UPDATES: 01 JUL 99 HIGHEST RN 226917-44-8

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 13, 1999

Please note that search-term pricing does apply when conducting SmartSELECT searches.

=> d stat que 120 L12 ST



Ak—N Ak—N—Ak Ak Ak Ak Ak Ak—N—Ak (21 22 (23 24 25) | 31 Ak—N—Ak (26 27 28 (230) 32 Ak 34

VAR G1=CH2/7/10
VAR G2=N/14/17/21/23/26/30
NODE ATTRIBUTES:
CONNECT IS M1 RC AT 1
CONNECT IS M1 RC AT 3
CONNECT IS M1 RC AT 5
CONNECT IS M1 RC AT 6
DEFAULT MLEVEL IS ATOM
DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RSPEC 1

NUMBER OF NODES IS 33

STEREO ATTRIBUTES: NONE

L14 436477 SEA FILE=REGISTRY ABB=ON PLU=ON OC5/ES

L16 SCR 1992

L18 SCR 2016 OR 2026 OR 2021

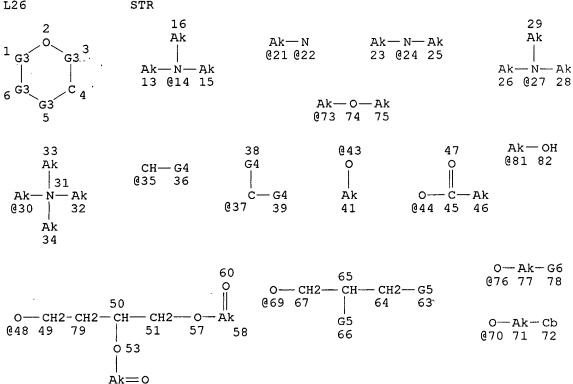
L20 60455 SEA FILE=REGISTRY SUB=L14 CSS FUL L12 AND L16 NOT L18

100.0% PROCESSED 148837 ITERATIONS

SEARCH TIME: 00.00.17

60455 ANSWERS

L26 HAS NO ANSWERS L26 STI



VAR G3=CH2/35/37 VAR G4=OH/43/AK/44/48/69/70/N/14/21/24/27/73/70/76/81/30 VAR G5=OH/43 VAR G6=N/22/24/27 NODE ATTRIBUTES: CONNECT IS M1 RC AT 4 CONNECT IS M1 RC AT 72 DEFAULT MLEVEL IS ATOM GGCAT IS PCY AT 72

GRAPH ATTRIBUTES:

RSPEC 1

NUMBER OF NODES IS 63

DEFAULT ECLEVEL IS LIMITED

STEREO ATTRIBUTES: NONE

=> s 126 sub=120 css sam STRUCTURE TOO LARGE - SEARCH ENDED

54 56

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=> fil reg FILE 'REGISTRY' ENTERED AT 08:36:45 ON 01 JUL 1999 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 1999 American Chemical Society (ACS)

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TSCA INFORMATION NOW CURRENT THROUGH JANUARY 13, 1999

Please note that search-term pricing does apply when conducting SmartSELECT searches.

=> d ide can 17

ANSWER 1 OF 1 REGISTRY COPYRIGHT 1999 ACS L7

221247-53-6 REGISTRY RN

.beta.-D-Glucopyranoside, (9Z)-9-octadecenyl 2-amino-2-deoxy- (9CI) CN INDEX NAME)

FS STEREOSEARCH

MF C24 H47 N O5

SR

CA, CAPLUS, TOXLIT LC STN Files:

Absolute stereochemistry. Double bond geometry as shown.

OH HO OH

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

1: 130:242316 REFERENCE

=> d his 17-

(FILE 'REGISTRY' ENTERED AT 08:30:59 ON 01 JUL 1999)

1 S L6 AND C24H47NO5 L7

E C46H85NO9/MF

L8 4 S E3

FILE 'HCAOLD' ENTERED AT 08:36:28 ON 01 JUL 1999

0 S L7 L9

FILE 'HCAPLUS' ENTERED AT 08:36:31 ON 01 JUL 1999

L10 1 S L7

FILE 'USPATFULL' ENTERED AT 08:36:33 ON 01 JUL 1999

- page 36 claim 8

claim 7 claim 7 - cmpd NOT found

L11 0 S L7

FILE 'REGISTRY' ENTERED AT 08:36:45 ON 01 JUL 1999

=> fil hcaplus FILE 'HCAPLUS' ENTERED AT 08:36:54 ON 01 JUL 1999 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 1999 AMERICAN CHEMICAL SOCIETY (ACS)

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FILE COVERS 1967 - 1 Jul 1999 VOL 131 ISS 1 FILE LAST UPDATED: 1 Jul 1999 (19990701/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

This file supports REG1stRY for direct browsing and searching of all substance data from the REGISTRY file. Enter HELP FIRST for more information.

=> d all 110

```
L10 ANSWER 1 OF 1 HCAPLUS COPYRIGHT 1999 ACS
```

- AN 1999:194158 HCAPLUS
- DN 130:242316
- TI Hydrophobic glycosylamine derivatives, compositions, and methods for their use
- IN Mumper, Russell J.; Tagliaferri, Frank
- PA Genemedicine, Inc., USA
- SO PCT Int. Appl., 88 pp.

CODEN: PIXXD2

- DT Patent
- LA English
- IC ICM C07H015-00
- CC 63-6 (Pharmaceuticals)

Section cross-reference(s): 33

FAN.CNT 1

	PATENT NO.				KIND		DATE		APPLICATION NO. DATE										
D.T.	WO 9912945						10000010		770 00 771 0000 1000000										
ΡI				AZ		19990318		. WO 98-US18888 19980908											
		W:	AL,	AM,	ΑT,	AU,	ΑZ,	ВA,	BB,	BG,	BR,	BY,	CA,	CH,	CN,	CU,	CZ,	DE,	
			DK,	EE,	ES,	FI,	GB,	GE,	GH,	GM,	HR,	ΗU,	ID,	IL,	IS,	JP,	KE,	KG,	
			ΚP,	KR,	ΚZ,	LC,	LK,	LR,	LS,	LT,	LU,	LV,	MD,	MG,	MK,	MN,	MW,	MX,	
			NO,	NZ,	PL,	PT,	RO,	RU,	SD,	SE,	SG,	SI,	SK,	SL,	TJ,	TM,	TR,	TT,	
			UΆ,	υG,	US,	UZ,	VN,	YU,	ZW,	AM,	ΑZ,	BY,	KG,	ΚZ,	MD,	RU,	ТJ,	TM	
		RW:					MW,			-			-						
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			CM,	GΑ,	GN,	GW,	ML,	MR,	ΝE,	SN,	TD,	ΤG							
	AU 9893839				A	Al 19990329				AU 98-93839					19980908				

PRAI US 97-58259 19970908

WO 98-US18888 19980908

OS MARPAT 130:242316

```
AR
     The invention relates in part to hydrophobic glycosylamine derivs.,
     methods for synthesizing hydrophobic derivs., compns. comprising these
     derivs., and methods for delivering macromols., such as proteins,
     peptides, lipids, carbohydrates, peptidomimetics, org. mols., and nucleic
     acids, to cells by administering these compns. The compds., compns., and
     methods of the invention are particularly useful for gene therapy and
     cancer treatment. Compns. contq. 1-mono-oleyl-.beta.-D-glucosamine or
     1-monopalmityl-.beta.-D-glucosamine, plasmid DNA comprising an IL-2 gene,
     and DOPE reduced the growth rate in mice by 30% after 9 days and by 25%
     after 13 days, resp.
     glycosylamine macromol delivery gene therapy antitumor
ST
     Nucleotides, biological studies
ΙT
     RL: THU (Therapeutic use); BIOL (Biological study); USES (Uses)
        (analogs; compns. and synthesis of hydrophobic glycosylamine derivs.
        for delivery of macromol. compds. to cells)
IT
     Macromolecular compounds
     RL: THU (Therapeutic use); BIOL (Biological study); USES (Uses)
        (anionic; compns. and synthesis of hydrophobic glycosylamine derivs.
        for delivery of macromol. compds. to cells)
ΙT
     Antitumor agents
     Drug delivery systems
     Gene therapy
     Infusions (drug delivery systems)
     Inhalants (drug delivery systems)
     Injections (drug delivery systems)
     Intravenous injections
     Liposomes (drug delivery systems)
     Oral drug delivery systems
     Plasmids
     Reducing agents
        (compns. and synthesis of hydrophobic glycosylamine derivs. for
        delivery of macromol. compds. to cells)
ΙT
     DNA
     Nucleic acids
     Polynucleotides
     RL: THU (Therapeutic use); BIOL (Biological study); USES (Uses)
        (compns. and synthesis of hydrophobic glycosylamine derivs. for
        delivery of macromol. compds. to cells)
IΤ
     Cryoprotectants
        (formulations contg.; compns. and synthesis of hydrophobic
        glycosylamine derivs. for delivery of macromol. compds. to cells)
IT
     Interleukin 2
     RL: BSU (Biological study, unclassified); BIOL (Biological study)
        (gene encoding; compns. and synthesis of hydrophobic glycosylamine
        derivs. for delivery of macromol. compds. to cells)
IT
     7440-05-3, Palladium, uses
     RL: CAT (Catalyst use); USES (Uses)
        (catalyst; compns. and synthesis of hydrophobic glycosylamine derivs.
        for delivery of macromol. compds. to cells)
ΤТ
     221247-53-6P
                    221247-55-8P
     RL: PRP (Properties); SPN (Synthetic preparation); THU (Therapeutic use);
     BIOL (Biological study); PREP (Preparation); USES (Uses)
        (compns. and synthesis of hydrophobic glycosylamine derivs. for
        delivery of macromol. compds. to cells)
                                                         63000-69-1
                               36653-82-4, Hexadecanol
IT
     143-28-2, Oleyl alcohol
     138395-62-7
     RL: RCT (Reactant)
        (compns. and synthesis of hydrophobic glycosylamine derivs. for
```

delivery of macromol. compds. to cells)

IT 221247-52-5P 221247-54-7P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation) (compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

IT 9003-39-8, PVP

RL: THU (Therapeutic use); BIOL (Biological study); USES (Uses) (formulations contg., as cryoprotectant; compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

IT 57-88-5, Cholesterol, biological studies 2462-63-7, DOPE
RL: THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(formulations contg.; compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

IT 302-01-2, Hydrazine, reactions 1333-74-0, Hydrogen, reactions 16940-66-2, Sodium borohydrate

RL: RCT (Reactant)

(reducing agent; compns. and synthesis of hydrophobic glycosylamine derivs. for delivery of macromol. compds. to cells)

=> fil reg

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TSCA INFORMATION NOW CURRENT THROUGH JANUARY 13, 1999

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=> d sca 18

L8 .4 ANSWERS REGISTRY COPYRIGHT 1999 ACS

IN Butanoic acid, 4-[D-glucopyranosyl(1-oxooctadecyl)amino]-4-oxo-,
10-octadecenyl ester (9CI)

MF C46 H85 N O9

Absolute stereochemistry. Double bond geometry unknown.

" free "view of L8 for mol. form., rage 84, claim 7 L8 4 ANSWERS REGISTRY COPYRIGHT 1999 ACS

IN L-Glutamic acid, N-D-ribonoyl-, di-(9Z)-9-octadecenyl ester (9CI)

MF C46 H85 N O9

CI COM

PAGE 1-A

PAGE 1-B

$$=$$
 CH $-$ (CH $_2$) $_7-$ Me

L8 4 ANSWERS REGISTRY COPYRIGHT 1999 ACS

IN 13-Docosenamide, N-[1-[(.beta.-D-glucopyranosyloxy)methyl]-2-hydroxy-3,7heptadecadienyl]-2-hydroxy- (9CI)

MF C46 H85 N O9

Absolute stereochemistry. Double bond geometry unknown.

L8 4 ANSWERS REGISTRY COPYRIGHT 1999 ACS

IN Heneicosanamide, N-[(1S, 2R, 3E, 7E, 9E)-1-[(.beta.-D-glucopyranosyloxy)methyl]-2-hydroxy-8-methyl-3,7,9-heptadecatrienyl]-2-

hydroxy-, (2R)- (9CI) MF **C46 H85 N 09**

HO_

PAGE 1-B

$$-$$
 (CH₂)₁₈-Me

ALL ANSWERS HAVE BEEN SCANNED

=> fil reg

FILE 'REGISTRY' ENTERED AT 09:24:55 ON 01 JUL 1999
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STRUCTURE FILE UPDATES: 01 JUL 99 HIGHEST RN 226917-44-8 DICTIONARY FILE UPDATES: 01 JUL 99 HIGHEST RN 226917-44-8

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 13, 1999

Please note that search-term pricing does apply when conducting SmartSELECT searches.

=> d ide can tot 128

L28 ANSWER 1 OF 5 REGISTRY COPYRIGHT 1999 ACS

RN 221247-55-8 REGISTRY

CN .beta.-D-Glucopyranoside, hexadecyl 2-amino-2-deoxy- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C22 H45 N O5

SR CA

LC STN Files: CA, CAPLUS, TOXLIT

Absolute stereochemistry.

from refore 2

.

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 130:242316

L28 ANSWER 2 OF 5 REGISTRY COPYRIGHT 1999 ACS

RN 221247-54-7 REGISTRY

CN .beta.-D-Glucopyranoside, hexadecyl 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C36 H53 N O10

SR CA

LC STN Files: CA, CAPLUS, TOXLIT

Absolute stereochemistry.

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 130:242316

L28 ANSWER 3 OF 5 REGISTRY COPYRIGHT 1999 ACS

RN **221247-52-5** REGISTRY

CN .beta.-D-Glucopyranoside, (9Z)-9-octadecenyl 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C38 H55 N O10

SR CA

LC STN Files: CA, CAPLUS, TOXLIT

Absolute stereochemistry.

Double bond geometry as shown.

1 REFERENCES IN FILE CA (1967 TO DATE)

1 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 130:242316

L28 ANSWER 4 OF 5 REGISTRY COPYRIGHT 1999 ACS

RN 138395-62-7 REGISTRY

CN .alpha.-D-Glucopyranoside, methyl 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C15 H17 N O7

SR CA

LC STN Files: BEILSTEIN*, CA, CAPLUS, CASREACT, TOXLIT (*File contains numerically searchable property data)

Absolute stereochemistry.

2 REFERENCES IN FILE CA (1967 TO DATE)

2 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 130:242316

REFERENCE 2: 116:59828

L28 ANSWER 5 OF 5 REGISTRY COPYRIGHT 1999 ACS

RN 63000-69-1 REGISTRY

CN .alpha.-D-Glucopyranosyl bromide, 2-deoxy-2-(1,3-dihydro-1,3-dioxo-2H-isoindol-2-yl)-, 3,4,6-triacetate (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C20 H20 Br N O9

LC STN Files: BEILSTEIN*, CA, CAPLUS, CASREACT, CHEMINFORMRX, TOXLIT, USPATFULL

(*File contains numerically searchable property data)

Absolute stereochemistry.

23 REFERENCES IN FILE CA (1967 TO DATE)

23 REFERENCES IN FILE CAPLUS (1967 TO DATE)

REFERENCE 1: 130:242316

REFERENCE 2: 129:343650

REFERENCE 3: 126:47484

REFERENCE 4: 126:8452

REFERENCE 5: 125:301540

REFERENCE 6: 125:196166

REFERENCE 7: 125:87030

REFERENCE 8: 123:33556

REFERENCE 9: 122:106316

REFERENCE 10: 121:80970

FILE 'HCAOLD' ENTERED AT 16:38:41 ON 01 JUL 1999 L20 137 S L19 FILE 'HCAPLUS' ENTERED AT 16:39:26 ON 01 JUL 1999 L21 1061 S L19 178 S L21 AND P/DT L22 L23 158 S L22 AND PY<1998 L24 77 S L22 AND US/PC L25 73 S L23 AND L24 77 S L24 AND (AY<1998 OR AY.B<1998 OR PRY<1998 OR PRY.B<1998) L26 77 S L25, L26 L27

FILE 'REGISTRY' ENTERED AT 16:40:55 ON 01 JUL 1999

=> fil hcaold

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PRE-1967 CHEMICAL ABSTRACTS FILE WITH HOUR-BASED PRICING FILE COVERS 1907-1966 FILE LAST UPDATED: 01 May 1997 (19970501/UP)

This file contains CAS Registry Numbers for easy and accurate substance identification. Title keywords, authors, and patent assignees are now searchable from 1907-1966. TIFF images of CA abstracts printed between 1907-1966 are available in the PAGE display formats.

This file supports REG1stRY for direct browsing and searching of all substance data from the REGISTRY file. Enter HELP FIRST for more information.

=> d 120 1-10 all hitstr

```
L20 ANSWER 1 OF 137 COPYRIGHT 1999 ACS
    CA65:20203f CAOLD
AN
     synthesis and antimetabolic properties of 4-N-hydroxycytosine derivs.
ΤI
ΑU
    Nelson, Donald J.
     antibiotics and related substances - (XXIV) amino sugars (11)
TΙ
     configurational studies of amino sugar glycosides and aminocyclitols by a
     Cu complex method
    Umezawa, Sumio; Tsuchiya, T.; Tatsuta, K.
ΑIJ
                3055-46-7 3867-93-4 13374-24-8
IT
       59-01-8
     13374-25-9 13374-26-0 13374-27-1 13374-28-2
     13374-32-8 13374-33-9 13374-34-0 13374-35-1 14187-81-6 14684-99-2
     89584-01-0
     3867-93-4 13374-24-8 13374-25-9
IΤ
     13374-26-0
     3867-93-4 HCAOLD
RN
     .beta.-D-Glucopyranoside, methyl 2-amino-2-deoxy-, hydrochloride (9CI)
CN
     (CA INDEX NAME)
```

● HCl

RN 13374-24-8 HCAOLD

CN Glucopyranoside, methyl 2-amino-2-deoxy-4,6-di-O-methyl-, hydrochloride, .beta.-D- (8CI) (CA INDEX NAME)

Absolute stereochemistry.

● HCl

RN 13374-25-9 HCAOLD

CN Glucopyranoside, methyl 2-amino-2-deoxy-6-O-methyl-, hydrochloride, .beta.-D- (8CI) (CA INDEX NAME)

Absolute stereochemistry.

HCl

RN 13374-26-0 HCAOLD

CN Glucopyranoside, methyl 2-amino-2-deoxy-3-0-methyl-, hydrochloride,

.beta.-D- (8CI) (CA INDEX NAME)

Absolute stereochemistry.

HCl

IT **90-77-7**RN 90-77-7 HCAOLD

CN D-Glucopyranose, 2-amino-2-deoxy- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

```
L20 ANSWER 3 OF 137 COPYRIGHT 1999 ACS
AN CA65:15476f CAOLD
TI synthesis and chemistry of 4-amino-4
```

TI synthesis and chemistry of 4-amino-4,6-dideoxy sugars - (II) glucose AU Stevens, Calvin L.; Blumbergs, P.; Daniher, F. A.; Otterbach, D. H.; Taylor, K. G.

6988-42-7 13042-62-1 **13042-64-3** 13042-65-4 13225-23-5 TΨ 13225-30-4 **13225-31-5 13225-32-6** 13231-14-6 13231-21-5 13231-22-6 13231-15-7 13231-16-8 13231-17-9 13231-20-4 13231-28-2 13231-29-3 13231-30-6 13231-23-7 13231-24-8 13231-27-1 13231-35-1 13231-36-2 13231-37-3 13231-31-7 13231-33-9 13231-34-0 13231-42-0 13231-43-1 13231-38-4 13231-39-5 13231-40-8 13405-51-1 13405-52-2 13518-22-4 14532-80-0 14532-81-1

IT 13042-64-3 13225-31-5 13225-32-6

RN 13042-64-3 HCAOLD

CN .alpha.-D-Glucopyranoside, methyl 4-amino-4,6-dideoxy- (9CI) (CA INDEX

NAME)

Absolute stereochemistry.

RN 13225-31-5 HCAOLD

CN Galactopyranoside, methyl 4-amino-4,6-dideoxy-, .alpha.-D- (8CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 13225-32-6 HCAOLD

CN Galactopyranoside, methyl 4-amino-4,6-dideoxy-, hydrochloride, .alpha.-D-(8CI) (CA INDEX NAME)

Absolute stereochemistry.

● HCl

L20 ANSWER 4 OF 137 COPYRIGHT 1999 ACS

AN CA65:12139c CAOLD

TI sulfonamides

PA Farbenfabriken Bayer A.-G.

DT Patent

```
ΤI
     sulfonamides
ΑU
     Horstmann, Harald; Wollweber, H.; Meng, K.
DT
     Patent
IT
     2439-57-8
                 4795-29-3
                              4795-30-6
                                          5064-46-0
                                                       6628-83-7
     7175-80-6
                 7175-81-7
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     14748-17-5
                 14796-39-5
                              71680-60-9
                              7179-99-9
ΙT
     6628-83-7
                 7179-97-7
     6628-83-7 HCAOLD
RN
CN
     2H-Pyran-2-methanamine, tetrahydro- (9CI) (CA INDEX NAME)
```

RN 7179-97-7 HCAOLD
CN 2H-Pyran-3-methanamine, tetrahydro-N-methyl- (9CI) (CA INDEX NAME)

RN 7179-99-9 HCAOLD
CN 2H-Pyran-3-methanamine, tetrahydro- (9CI) (CA INDEX NAME)

```
L20 ANSWER 5 OF 137 COPYRIGHT 1999 ACS
AN
     CA65:9369e CAOLD
    mannosamine in the cell wall of Clostridium welchii (Type A)
ΤI
ΑU
     Pickering, Brian T.
IT
     2636-92-2
IT
     2636-92-2
RN
     2636-92-2 HCAOLD
L20
    ANSWER 6 OF 137 COPYRIGHT 1999 ACS
ΑN
     CA65:7258g CAOLD
     amino sugars - (X) prepn. of 3-amino-3,6-dideoxy-D-gulose, -D-allose, and
TI
     -L-allose derivs.
ΑU
     Capek, K.; Jary, J.
                            7646-69-7 13042-61-0 13042-62-1 13042-63-2
     4613-56-3
                 4782-93-8
IT
     13042-64-3 13042-65-4 13042-66-5 13042-67-6 13042-68-7
     13042-69-8 13042-70-1 13042-71-2 13042-72-3 13042-73-4
```

IT 13042-64-3

RN 13042-64-3 HCAOLD

CN .alpha.-D-Glucopyranoside, methyl 4-amino-4,6-dideoxy- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

L20 ANSWER 7 OF 137 COPYRIGHT 1999 ACS

AN CA65:3944f CAOLD

aminohexoses - (IX) N-deacetylation of methyl 2-acetamido-2-deoxy-.beta.-D-glucopyranoside, methyl 2-acetamido-2-deoxy-3-O-methyl-.alpha. (and .beta.)-D-glucopyranoside, and methyl 2-acetamido-2-deoxy-4,6-O-ethylidene-.alpha.-D-glucopyranoside

AU Fujinaga, Mamoru; Matsushima, Y.

IT 3055-46-7 **3867-92-3 3867-93-4** 6619-04-1 10300-75-1 10300-79-5 10300-80-8 10368-82-8 10427-77-7 10427-79-9 22594-60-1

IT 3867-92-3 3867-93-4

RN 3867-92-3 HCAOLD

CN .beta.-D-Glucopyranoside, methyl 2-amino-2-deoxy- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 3867-93-4 HCAOLD

CN .beta.-D-Glucopyranoside, methyl 2-amino-2-deoxy-, hydrochloride (9CI) (CA INDEX NAME)

HCl

```
L20 ANSWER 8 OF 137 COPYRIGHT 1999 ACS
     CA64:19745g CAOLD
ΑN
     studies on kasugamycin - (V) structure of kasugamycin
ΤI
     Suhara, Yasuji; Maeda, K.; Umezawa, H.; Ohno, M.
ΑU
ΙT
     6980-16-1
                6980-17-2
                            6980-18-3
                                         6980-19-4
     6991-09-9
                21256-68-8
                            94892-82-7
ΙT
     6980-19-4
                 6991-09-9 21256-68-8
     94892-82-7
     6980-19-4 HCAOLD
RN
     D-arabino-Hexopyranoside, methyl 2-amino-4-(cyanoamino)-2,3,4,6-tetradeoxy-
CN
     , .beta.- (8CI) (CA INDEX NAME)
```

Absolute stereochemistry.

$$\begin{array}{c} \text{OMe} \\ \text{H}_2\text{N} \\ \text{S} \\ \text{R} \\ \text{N} \\ \text{N} \\ \text{N} \end{array}$$

RN 6991-09-9 HCAOLD

CN D-arabino-Hexopyranoside, methyl 2-amino-4-[(1-carboxyformimidoyl)amino]-2,3,4,6-tetradeoxy-, monohydrochloride, .beta.- (8CI) (CA INDEX NAME)

● HCl

RN 21256-68-8 HCAOLD

CN D-arabino-Hexopyranoside, methyl 2-amino-4-(cyanoamino)-2,3,4,6-tetradeoxy-(7CI, 8CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 94892-82-7 HCAOLD

CN D-arabino-Hexopyranoside, methyl 2-amino-4-[(1-carboxyformimidoyl)amino]-2,3,4,6-tetradeoxy-, hydrochloride (7CI) (CA INDEX NAME)

$$H_2N$$
 S
 R
 Me
 HO_2C
 NH

HCl

L20 ANSWER 9 OF 137 COPYRIGHT 1999 ACS ΑN CA64:18009a CAOLD ΤI spectrophotometric assay for hexosamines ΑU Galambos, John T.; Shapira, R. IT 90-76-6 579-33-9 2636-92-2 3416-24-8 579-33-9 2636-92-2 IT 90-76-6 90-76-6 HCAOLD RN D-Galactopyranose, 2-amino-2-deoxy- (9CI) (CA INDEX NAME) CN

Absolute stereochemistry.

RN 579-33-9 HCAOLD CN D-Mannopyranose, 2-amino-2-deoxy- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 2636-92-2 HCAOLD

```
L20 ANSWER 10 OF 137 COPYRIGHT 1999 ACS
AN
     CA64:17696f CAOLD
TI
     potential cytostats - (VI) configuration and derivs. of the
     1-deoxy-1-methylaminoalditols obtained from aldohexoses and aldopentoses
ΑU
     Dorn, Helmut; Welfle, H.; Liebig, R.
                                          7084-27-7
IT
     7057-02-5
                 7057-03-6
                              7057-04-7
                                                      7115-35-7
                              7115-40-4
     7115-38-0
                 7115-39-1
                                          7115-41-5
                                                      7115-42-6
     7115-43-7
                 7115-45-9
                              7115-46-0
                                          7115-47-1
                                                      7115-48-2
                                                                   7115-49-3
     7115-50-6
                 7115-51-7
                              7115-52-8
                                          7115-53-9
                                                      7115-54-0
                                                                   7115-55-1
                 7115-57-3
                              7115-58-4
     7115-56-2
                                          7115-59-5
                                                      7115-60-8
                                                                   7115-61-9
     7115-62-0
                 7115-63-1
                              7115-64-2
                                          7115-65-3
                                                      7139-59-5
                                                                   7185-78-6
     7185-79-7
                 7233-42-3
                              7297-32-7
                                         79253-93-3
IT
     7115-35-7
                 7115-39-1
RN
     7115-35-7
                HCAOLD
     Ribopyranose, 2-amino-2-deoxy-, .alpha.-D- (7CI, 8CI) (CA INDEX NAME)
CN
```

Absolute stereochemistry.

RN 7115-39-1 HCAOLD

CN Ribopyranose, 2-amino-2-deoxy-, hydrochloride, .alpha.-D- (7CI, 8CI) (CA INDEX NAME)

Absolute stereochemistry.

HCl